

Succinic acid, cyclohexylmethyl 2-methoxyethyl ester

Inchi:	InChI=1S/C14H24O5/c1-17-9-10-18-13(15)7-8-14(16)19-11-12-5-3-2-4-6-12/h12H,2-11H
InchiKey:	USYVFSYSGOZHKE-UHFFFAOYSA-N
Formula:	C14H24O5
SMILES:	COCCOC(=O)CCC(=O)OCC1CCCCC1
Mol. weight [g/mol]:	272.34

Physical Properties

Property code	Value	Unit	Source
gf	-481.39	kJ/mol	Joback Method
hf	-899.79	kJ/mol	Joback Method
hfus	30.61	kJ/mol	Joback Method
hvap	67.91	kJ/mol	Joback Method
log10ws	-2.15		Crippen Method
logp	2.080		Crippen Method
mvol	218.010	ml/mol	McGowan Method
pc	1920.30	kPa	Joback Method
rinpol	1967.00		NIST Webbook
rinpol	1967.00		NIST Webbook
tb	714.27	K	Joback Method
tc	912.84	K	Joback Method
tf	421.47	K	Joback Method
vc	0.819	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	645.13	J/molxK	714.27	Joback Method
cpg	662.57	J/molxK	747.36	Joback Method
cpg	678.94	J/molxK	780.46	Joback Method
cpg	694.24	J/molxK	813.55	Joback Method
cpg	708.46	J/molxK	846.65	Joback Method
cpg	721.61	J/molxK	879.74	Joback Method
cpg	733.67	J/molxK	912.84	Joback Method
dvisc	0.0012192	Paxs	421.47	Joback Method

dvisc	0.0006410	Paxs	470.27	Joback Method
dvisc	0.0003803	Paxs	519.07	Joback Method
dvisc	0.0002468	Paxs	567.87	Joback Method
dvisc	0.0001715	Paxs	616.67	Joback Method
dvisc	0.0001257	Paxs	665.47	Joback Method
dvisc	0.0000961	Paxs	714.27	Joback Method

Sources

McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U390742&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws
Joback Method:	https://en.wikipedia.org/wiki/Joback_method

Legend

cp_g:	Ideal gas heat capacity
dvisc:	Dynamic viscosity
g_f:	Standard Gibbs free energy of formation
h_f:	Enthalpy of formation at standard conditions
h_{fus}:	Enthalpy of fusion at standard conditions
h_{vap}:	Enthalpy of vaporization at standard conditions
log₁₀ws:	Log ₁₀ of Water solubility in mol/l
log_p:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume
pc:	Critical Pressure
rinpol:	Non-polar retention indices
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

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